

	Final Report
V	Revised Report

Report Date: 08-Feb-18 17:05

Laboratory Report SC43423

Gulf Oil L.P. 281 Eastern Avenue Chelsea, MA 02150 Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA

Project #: Gulf Chelsea

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87936 Maine # MA138 New Hampshire # 2972/2538 New Jersey # MA011 New York # 11393 Pennsylvania # 68-04426/68-02924 Rhode Island # LAO00348 USDA # P330-15-00375 Vermont # VT-11393



Authorized by:

Christina White Technical Director

Christina a. White

Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 13 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

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Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.

Sample Summary

Work Order: SC43423

Project: Gulf Terminal - Chelsea, MA

Project Number: Gulf Chelsea

Laboratory ID	Client Sample ID	<u>Matrix</u>	Date Sampled	Date Received
SC43423-01	Outfall 003	Surface Water	25-Jan-18 08:45	25-Jan-18 14:40
SC43423-02	TB-1 (Trip Blank)	Aqueous	25-Jan-18 00:00	25-Jan-18 14:40

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CASE NARRATIVE:

Data has been reported to the RDL. This report excludes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the reporting limit are reported as "<" (less than) the reporting limit in this report.

The samples were received 1.6 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 1.0 degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group. If method or program required MS/MSD/Dup were not performed, sufficient sample was not provided to the laboratory.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

February 8, 2018 Report Revision Case Narrative:

This report has been revised to update the analyte list for 8270 as requested.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

SW846 8260C

Calibration:

1801070

Analyte quantified by quadratic equation type calibration.

Naphthalene

This affected the following samples:

1801129-BLK1

1801129-BS1

1801129-BSD1

Outfall 003

S816062-ICV1

S816241-CCV1

TB-1 (Trip Blank)

SW846 8270D SIM

Samples:

S816305-CCV1

Analyte percent difference is outside individual acceptance criteria (20), but within overall method allowances.

Indeno (1,2,3-cd) pyrene (-25.2%)

This affected the following samples:

1801124-BLK1

1801124-BS1

1801124-BSD1

This laboratory report is not valid without an authorized signature on the cover page.

Sample Acceptance Check Form

Project:	Gulf Terminal - Chelsea, MA / Gulf Chelsea			
Work Order:	SC43423			
Sample(s) received on:	1/25/2018			
The following outlines th	ne condition of samples for the attached Chain of Custody upon receipt.			
		Yes	<u>No</u>	<u>N/A</u>
Were custody se	als present?		\checkmark	
Were custody se	als intact?			✓
Were samples re	ceived at a temperature of $\leq 6^{\circ}$ C?	\checkmark		
Were samples re	frigerated upon transfer to laboratory representative?	\checkmark		
Were sample con	ntainers received intact?	\checkmark		
	operly labeled (labels affixed to sample containers and include sample ID, site project number and the collection date)?	$\overline{\checkmark}$		
Were samples ac	companied by a Chain of Custody document?	\checkmark		
include sample I	ustody document include proper, full, and complete documentation, which shall D, site location, and/or project number, date and time of collection, collector's name, e, sample matrix and any special remarks concerning the sample?			
Did sample cont	ainer labels agree with Chain of Custody document?	\checkmark		
Were samples re	ceived within method-specific holding times?	\overline{V}	П	

Client:

Gulf Oil L.P.

Summary of Hits

Client ID:

Outfall 003

Lab ID: SC43423-01

Flag Parameter Result **Reporting Limit** Units **Analytical Method** Total Suspended Solids 25.8 1.2 SM2540D (11) mg/l 1.2 1.0 SW846 8260C Benzene $\mu g/l$ 0.198 Naphthalene 0.047 $\mu g/l$ SW846 8270D SIM

Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.

Sample Id Outfall 00 SC43423-				Client P Gulf C	-		Matrix Surface W		ection Date 5-Jan-18 08			<u>cceived</u> -Jan-18	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Volatile O	rganic Compounds												
Volatile O	rganic Aromatics by SW8	46 8260											
Prepared	by method SW846 5030	Water MS											
71-43-2	Benzene	1.2		μg/l	1.0	0.3	1	SW846 8260C	26-Jan-18	30-Jan-18	GMA	1801129	
91-20-3	Naphthalene	< 1.0		μg/l	1.0	0.4	1	II .	"	"	"	"	
Surrogate	recoveries:												
460-00-4	4-Bromofluorobenzene	111			70-13	80 %		"	"	"	"	"	
2037-26-5	Toluene-d8	102			70-13	80 %		n	"	"		"	
17060-07-0	1,2-Dichloroethane-d4	99			70-13	80 %		"	u	"	"	"	
1868-53-7	Dibromofluoromethane	105			70-13	80 %		n .	"	"	"	"	
Semivolati	ile Organic Compounds by	GCMS											
SVOCs b	y SIM												
Prepared	by method SW846 35100	<u>2</u>											
50-32-8	Benzo (a) pyrene	< 0.047		μg/l	0.047	0.019	1	SW846 8270D SIM	26-Jan-18	26-Jan-18	MSL	1801124	
91-20-3	Naphthalene	0.198		μg/l	0.047	0.020	1	n .	"	"	"	"	
Surrogate	recoveries:												
205440-82-0	Benzo (e) pyrene-d12	72			30-13	80 %		II .	u u	"	"	"	
	le Petroleum Hydrocarbon by method General Prepa		<u> </u>										
	Oil & Grease	< 1.00	OG	mg/l	1.00	0.915	1	EPA 1664B	05-Feb-18	06-Feb-18	DJS	1801626	Х
General C	hemistry Parameters												
	рH	6.63	рН	pH Units			1	ASTM D 1293-99B	25-Jan-18 17:00	25-Jan-18 17:30	BD	1801116	Χ
	Total Suspended Solids	25.8		mg/l	1.2	0.5	1	SM2540D (11)	26-Jan-18	30-Jan-18	СМВ	1801137	Х

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Sample Id TB-1 (Tri SC43423-	•				Project # Chelsea		Matrix Aqueou		ection Date -Jan-18 00			<u>ceived</u> Jan-18	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Volatile Or	rganic Compounds												
	rganic Aromatics by SW8 by method SW846 5030												
71-43-2	Benzene	< 1.0		μg/l	1.0	0.3	1	SW846 8260C	26-Jan-18	30-Jan-18	GMA	1801129	
91-20-3	Naphthalene	< 1.0		μg/l	1.0	0.4	1	п	"	"	"	u	
Surrogate r	recoveries:												
460-00-4	4-Bromofluorobenzene	107			70-13	0 %		"	"	"	"	"	
2037-26-5	Toluene-d8	101			70-13	0 %		"	"	"		"	
17060-07-0	1,2-Dichloroethane-d4	99			70-13	0 %		"	"	"	"	"	
1868-53-7	Dibromofluoromethane	106			70-13	0 %		п	"	"	u	u	

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Volatile Organic Compounds - Quality Control

nalyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limi
W846 8260C										
atch 1801129 - SW846 5030 Water MS										
Blank (1801129-BLK1)					Pre	epared: 26-	Jan-18 An	alyzed: 29-Ja	<u>an-18</u>	
Benzene	< 1.0		μg/l	1.0						
Naphthalene	< 1.0		μg/l	1.0						
Surrogate: 4-Bromofluorobenzene	54.6		μg/l		50.0		109	70-130		
Surrogate: Toluene-d8	50.2		μg/l		50.0		100	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.4		μg/l		50.0		99	70-130		
Surrogate: Dibromofluoromethane	53.2		μg/l		50.0		106	70-130		
LCS (1801129-BS1)					Pre	epared: 26-	Jan-18 An	alyzed: 29-Ja	<u>ın-18</u>	
Benzene	20.7		μg/l		20.0		103	70-130		
Naphthalene	20.0		μg/l		20.0		100	70-130		
Surrogate: 4-Bromofluorobenzene	52.5		μg/l		50.0		105	70-130		
Surrogate: Toluene-d8	50.8		μg/l		50.0		102	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.6		μg/l		50.0		99	70-130		
Surrogate: Dibromofluoromethane	53.2		μg/l		50.0		106	70-130		
LCS Dup (1801129-BSD1)					Pre	epared: 26-	Jan-18 An	alyzed: 29-Ja	n-18	
Benzene	19.4		μg/l		20.0		97	70-130	7	20
Naphthalene	18.1		μg/l		20.0		91	70-130	10	20
Surrogate: 4-Bromofluorobenzene	51.8		μg/l		50.0		104	70-130		
Surrogate: Toluene-d8	50.3		μg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	50.4		μg/l		50.0		101	70-130		
Surrogate: Dibromofluoromethane	52.3		μg/l		50.0		105	70-130		

Semivolatile Organic Compounds by GCMS - Quality Control

nalyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
W846 8270D SIM										
atch 1801124 - SW846 3510C										
Blank (1801124-BLK1)					Pre	epared & A	nalyzed: 26-	Jan-18		
Acenaphthene	< 0.051		μg/l	0.051						
Acenaphthylene	< 0.051		μg/l	0.051						
1-Methylnaphthalene	< 0.051		μg/l	0.051						
Anthracene	< 0.051		μg/l	0.051						
Benzo (a) anthracene	< 0.051		μg/l	0.051						
Benzo (a) pyrene	< 0.051		μg/l	0.051						
Benzo (b) fluoranthene	< 0.051		μg/l	0.051						
Benzo (g,h,i) perylene	< 0.051		μg/l	0.051						
Benzo (k) fluoranthene	< 0.051		μg/l	0.051						
Chrysene	< 0.051		μg/l	0.051						
Dibenzo (a,h) anthracene	< 0.051		μg/l	0.051						
Fluoranthene	< 0.051		μg/l	0.051						
Fluorene	< 0.051		μg/l	0.051						
Indeno (1,2,3-cd) pyrene	< 0.051		μg/l	0.051						
2-Methylnaphthalene	< 0.051		μg/l	0.051						
Naphthalene	< 0.051		μg/l	0.051						
Phenanthrene	< 0.051		μg/l	0.051						
Pyrene	< 0.051		μg/l	0.051						
Surrogate: Benzo (e) pyrene-d12	0.687		μg/l		1.01		68	30-130		
LCS (1801124-BS1)			1.0			epared & A	nalyzed: 26-			
Acenaphthene	0.535		μg/l	0.051	1.01	•	53	40-140		
Acenaphthylene	0.551		μg/l	0.051	1.01		55	40-140		
1-Methylnaphthalene	0.539		μg/l	0.051	1.01		53	40-140		
Anthracene	0.724		μg/l	0.051	1.01		72	40-140		
Benzo (a) anthracene	0.897		μg/l	0.051	1.01		89	40-140		
Benzo (a) pyrene	0.716		μg/l	0.051	1.01		71	40-140		
Benzo (b) fluoranthene	0.752		μg/l	0.051	1.01		74	40-140		
Benzo (g,h,i) perylene	0.635		μg/l	0.051	1.01		63	40-140		
Benzo (k) fluoranthene	0.822		μg/l	0.051	1.01		81	40-140		
Chrysene	0.787		μg/l	0.051	1.01		78	40-140		
Dibenzo (a,h) anthracene	0.700		μg/l	0.051	1.01		69	40-140		
Fluoranthene	0.820		μg/l	0.051	1.01		81	40-140		
Fluorene	0.737		μg/l	0.051	1.01		73	40-140		
Indeno (1,2,3-cd) pyrene	0.702		μg/l	0.051	1.01		70	40-140		
2-Methylnaphthalene	0.659		μg/l	0.051	1.01		65	40-140		
Naphthalene	0.506		μg/l	0.051	1.01		50	40-140		
Phenanthrene	0.836		μg/l	0.051	1.01		83	40-140		
Pyrene	0.826		μg/l	0.051	1.01		82	40-140		
Surrogate: Benzo (e) pyrene-d12	0.899		μg/l		1.01		89	30-130		
LCS Dup (1801124-BSD1)			P3.1			epared & A	nalyzed: 26-			
Acenaphthene	0.531		μg/l	0.050	1.00		53	40-140	0.8	20
Acenaphthylene	0.548		μg/l	0.050	1.00		55	40-140	0.5	20
1-Methylnaphthalene	0.522		μg/l	0.050	1.00		52	40-140	3	20
Anthracene	0.684		μg/l	0.050	1.00		68	40-140	6	20
Benzo (a) anthracene	0.815		μg/l	0.050	1.00		82	40-140	10	20
Benzo (a) pyrene	0.698		μg/l	0.050	1.00		70	40-140	3	20
Benzo (b) fluoranthene	0.772		μg/l	0.050	1.00		77	40-140	3	20
Benzo (g,h,i) perylene	0.671		μg/l	0.050	1.00		67	40-140	5	20
Benzo (k) fluoranthene	0.820		μg/l	0.050	1.00		82	40-140	0.3	20
Chrysene	0.820		μg/l μg/l	0.050	1.00		82	40-140	4	20

Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Analyte(s)	Result	Tag	Units	KDL	Level	Result	/0KEC	Lillits	KLD	Lillit
SW846 8270D SIM										
Batch 1801124 - SW846 3510C										
LCS Dup (1801124-BSD1)					Pro	epared & A	nalyzed: 26	-Jan-18		
Dibenzo (a,h) anthracene	0.755		μg/l	0.050	1.00		76	40-140	8	20
Fluoranthene	0.763		μg/l	0.050	1.00		76	40-140	7	20
Fluorene	0.730		μg/l	0.050	1.00		73	40-140	1	20
Indeno (1,2,3-cd) pyrene	0.707		μg/l	0.050	1.00		71	40-140	0.7	20
2-Methylnaphthalene	0.588		μg/l	0.050	1.00		59	40-140	11	20
Naphthalene	0.509		μg/l	0.050	1.00		51	40-140	0.6	20
Phenanthrene	0.835		μg/l	0.050	1.00		84	40-140	0.2	20
Pyrene	0.836		μg/l	0.050	1.00		84	40-140	1	20
Surrogate: Benzo (e) pyrene-d12	0.850		μg/l		1.00		85	30-130		

Extractable Petroleum Hydrocarbons - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
EPA 1664B										
Batch 1801626 - General Preparation SVOC										
Blank (1801626-BLK1)					Pre	epared: 05-F	eb-18 An	alyzed: 06-F	eb-18	
Oil & Grease	< 1.03		mg/l	1.03						
LCS (1801626-BS1)					Pre	epared: 05-F	eb-18 An	alyzed: 06-F	eb-18	
Oil & Grease	36.1		mg/l	1.02	40.4		89	78-114		

General Chemistry Parameters - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
ASTM D 1293-99B										
Batch 1801116 - General Preparation										
Reference (1801116-SRM1)					Pre	epared & Ar	nalyzed: 25	-Jan-18		
рН	6.05		pH Units		6.00		101	97.5-102. 5		
Reference (1801116-SRM2)					Pre	epared & Ar	nalyzed: 25	-Jan-18		
рН	6.01		pH Units		6.00		100	97.5-102. 5		
SM2540D (11)										
Batch 1801137 - General Preparation										
Blank (1801137-BLK1)					Pre	epared: 26-	Jan-18 An	alyzed: 30-Ja	<u>ın-18</u>	
Total Suspended Solids	< 0.5		mg/l	0.5						
LCS (1801137-BS1)					Pre	epared: 26-	Jan-18 An	alyzed: 30-Ja	<u>an-18</u>	
Total Suspended Solids	92.0		mg/l	10.0	100		92	90-110		

Notes and Definitions

dry Sample results reported on a dry weight basis

NR Not Reported

RPD Relative Percent Difference

OG The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed

when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample

volume was submitted to fulfill the requirement.

pH The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as

soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt.

All soil samples are analyzed as soon as possible after sample receipt.

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

<u>Matrix Spike</u>: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

<u>Continuing Calibration Verification:</u> The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.

eurofins	ofins	CHAIN	CHAIN OF CUSTODY RECORD		Special Handling: Standard TAT - 7 to 10 business days
	Spectrum Analytical) A	Page 1 of 1		All TAT's subject to laboratory approval Min. 24-hr notification needed for rushes Samples disposed after 30 days unless otherwise instructed.
Report To:	Andrew Adams	Invoice To:	Christopher Gill	Project No:	Gulf Chelsea
	Quit Oil LP		Colt Oil LP	Site Name:	Gulf Chelsea Termina)
	Chelsea MA 02150	1 1	Wellesley MA . 02481-3705	Location:	28) Eastern Luc. State: MA
Project Mgr:	Andrew Adams	P.O No.:	Quote #:	Sampler(s).	AECOM 1000 Marie 1

Please send report to	* Please se	tor	Corecction Factor	0	144	2	12					The But	J. J.	
enrifer atking accom. com	E-mail to: jennifer.		Observed	2	091	2	35	/	4	HIL R	1	//	1401	
	EDD format:		Temp °C	e:	Time:		Date:			Received by:		Relinquished by:	Rélinqui	
				2		- 1								
							ly c	120						
						576								
						BIG	uliz						6	
							بو				TripBlank	TB-1 (T		
and the	<		230	ty 8	est.	_		<	2180	1-2518	003	Outfull 003		
		<	46.0	100 j	-				5480	1-25-18	203	Outen 003		
1 mistery		1		44	-			1980	2480	1-25-18 6	003	Outful 003		
and broken			<			9			5480	1-25-18	003	Outfall 003		
- Mentana is SW	, ,			1		958	3	Ses	5480	1-25-18	003	Outfall 003	4342301	2
C Other: State-specific reporting standards:	(-	-	1			# of		Time:	Date:	9:	Sample ID:	Lab ID:	
-	5 5	TSS PH	PAI	OC:	Clear Plastic		VOA	/pe itrix		C=Compsite		G= Grab	G= 0	
			+5			r Glas	Vials			X3=	X2=		X1=	
· St				nz,	¥	S	pjej"		35	Air SG=Soil Gas	A=Indoor/Ambient Air	SL=Sludge A=	0=0il S0 =Soil	
Report? Yes	Analysis				Containers	Con			WW=Waste Water		r SW=Surface Water	GW=Groundwater	DW=Drinking Water	
* additional charges may appply	3			2								**		
QA/QC Reporting Notes:	List Preservative Code below:	List Prese					cid	6=Ascorbic Acid		4=HNO ₃ 5=NaOH	3=H ₂ SO ₄	1=Na ₂ S2O ₃	F=Field Filtered 1=Na ₂ S2O ₃ 2=HCl 3=H ₂ SO ₄ 7=CH3OH 8=NaHSO ₂ 9=Dejonized Water 10=H ₂ PO ₂	
	AECOM.				te #:	Quote #:		F 10	P.O No.:		Adams		Project Mgr:	
Jos Barragan	ampler(s).	_				,					こくもし	CXX カーン	Talanhana #	

Condition upon receipt: Custody Seals:

Present
Intact
Broken

Batch Summary

1801116

General Chemistry Parameters

1801116-SRM1 1801116-SRM2

SC43423-01 (Outfall 003)

1801124

Semivolatile Organic Compounds by GCMS

1801124-BLK1 1801124-BS1 1801124-BSD1

SC43423-01 (Outfall 003)

1801129

Volatile Organic Compounds

1801129-BLK1 1801129-BSD1 1801129-BSD1 SC43423-01 (Outfall 003) SC43423-02 (TB-1 (Trip Blank))

1801137

General Chemistry Parameters

1801137-BLK1 1801137-BS1 SC43423-01 (Outfall 003)

1801626

Extractable Petroleum Hydrocarbons

1801626-BLK1 1801626-BS1

SC43423-01 (Outfall 003)

S711062

Semivolatile Organic Compounds by GCMS

S711062-CAL1

S711062-CAL2

S711062-CAL3

S711062-CAL4

S711062-CAL5

S711062-CAL6

S711062-CAL7

S711062-CAL8

S711062-CAL9

S711062-ICV1

S711062-LCV1

S711062-LCV2

S711062-TUN1

S816062

Volatile Organic Compounds

S816062-CAL1

S816062-CAL2

S816062-CAL3

S816062-CAL4

S816062-CAL5

S816062-CAL6

S816062-CAL7

S816062-CAL8

S816062-CAL9

S816062-ICV1

S816062-LCV1

S816062-LCV2

S816062-TUN1

S816241

Volatile Organic Compounds

S816241-CCV1

S816241-TUN1

S816305

Semivolatile Organic Compounds by GCMS

S816305-CCV1

S816305-TUN1